

MARKED UP COPY OF AMENDMENT

IN THE SPECIFICATION

Page 7, delete line 11.

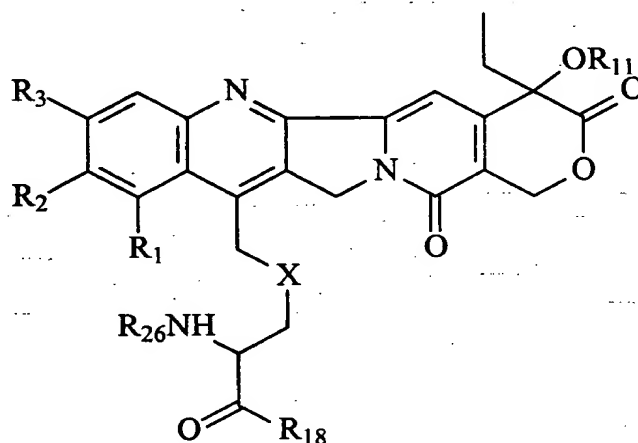
Page 12, delete line 8.

Page 19, delete line 3.

IN THE CLAIMS

Please amend the claims as follows:

--1. (Twice Amended) A compound comprising:



wherein R₁ and R₂, are each independently

NO₂, NH₂, H, F, Cl, Br, I, COOH, OH, O-C₁₋₆ alkyl, SH, S-C₁₋₆ alkyl, CN, NH-C₁₋₆ alkyl, N(C₁₋₆ alkyl)₂, CHO, C₁₋₈ alkyl, N₃,

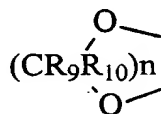
-Z-(CH₂)_a-N-((CH₂)_bOH)₂, wherein Z is selected from the group consisting of O, NH and S, and a and b are each independently an integer of 2 or 3,

-Z-(CH₂)_a-N-(C₁₋₆ alkyl)₂ wherein Z is selected from the group consisting of O, NH and S, and a is an integer of 2 or 3,

-CH₂NR₄R₅, where (a) R₄ and R₅ are, independently, hydrogen, C₁₋₆ alkyl, C₃₋₇ cycloalkyl, C₃₋₇ cycloalkyl-C₁₋₆ alkyl, C₂₋₆ alkenyl, hydroxy-C₁₋₆ alkyl, C₁₋₆ alkoxy-C₁₋₆ COR₆ where R₆ is hydrogen, C₁₋₆ alkyl, perhalo-C₁₋₆ alkyl, C₃₋₇ cycloalkyl, C₃₋₇ cycloalkyl-C₁₋₆ alkyl, C₂₋₆ alkenyl, hydroxy-C₁₋₆ alkyl, C₁₋₆ alkoxy, C₁₋₆ alkoxy-C₁₋₆ alkyl, or (b) R₄ and R₅ taken together with the nitrogen atom to which they are attached form a saturated 3-7 membered heterocyclic ring which may contain a O, S or NR₇ group, where R₇ is hydrogen, C₁₋₆ alkyl, perhalo-C₁₋₆ alkyl, aryl, aryl substituted with one or more groups selected from the group consisting of C₁₋₆ alkyl, halogen, nitro, amino, C₁₋₆ alkylamino, perhalo-C₁₋₆ alkyl, hydroxy-C₁₋₆ alkyl, C₁₋₆ alkoxy, C₁₋₆ alkoxy-C₁₋₆ alkyl and -COR₈ where R₈ is hydrogen, C₁₋₆ alkyl perhalo-C₁₋₆ alkyl, C₁₋₆ alkoxy, aryl, and aryl substituted with one or more C₁₋₆ alkyl, perhalo-C₁₋₆ alkyl, hydroxy-C₁₋₆ alkyl, or C₁₋₆ alkoxy-C₁₋₆ alkyl groups;

R₃ is H; or

or R₂ and R₃ combine to form a ring



where R₉ and R₁₀ are each independently H or F and n is an integer of 1 or 2;

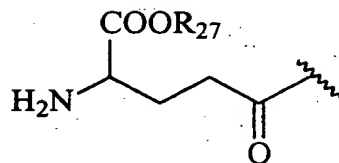
R₁₁ is H, or C(O)-(CH₂)_m-NR₁₂R₁₃, where m is an integer of 1-6 or

-C(O)CHR₁₄NR₁₂R₁₃, where R₁₄ is the side chain of one of the naturally occurring α-amino acids, R₁₂ and R₁₃ are, independently, hydrogen, C₁₋₈ alkyl or -C(O)CHR₁₅NR₁₆R₁₇, where R₁₅ is the side chain of one of the naturally occurring α-amino acids and R₁₆ and R₁₇ are each independently hydrogen or C₁₋₈ alkyl;

R₁₈ is OR₁₉ or R₁₉OC(O)-(CH₂)_m-NR₂₀, or R₂₁OC(O)CHR₂₂NR₂₀, where R₁₉ is H or C₁₋₆ alkyl, m is an integer of 1-6, R₂₂ is the side chain of one of the naturally occurring α-amino acids, R₂₀ is hydrogen, C₁₋₈ alkyl or -C(O)CHR₂₃NR₂₄R₂₅, where R₂₃ is the side chain of one

of the naturally occurring α -amino acids and R_{24} and R_{25} are each independently hydrogen or C_{1-8} alkyl;

R_{26} is H or



where R_{27} is H or C_{1-6} alkyl; and

X is S or O,

[provided that R_{18} and R_{26} are not both H;]

or a pharmaceutically acceptable salt thereof.--